=> d his

L7

(FILE 'HOME' ENTERED AT 18:23:40 ON 19 DEC 2005)

FILE 'REGISTRY' ENTERED AT 18:23:45 ON 19 DEC 2005

STRUCTURE UPLOADED L1

L2 2 S L1

8 S L1 FULL L3

FILE 'HCAPLUS' ENTERED AT 18:24:45 ON 19 DEC 2005

L4

FILE 'HCAPLUS, HCAOLD, USPATFULL, EPFULL, BIOSIS, MEDLINE' ENTERED AT 18:26:01 ON 19 DEC 2005

L5 913 S ?HYPERFORIN OR ?ADHYPERFORIN

'94 S L5 AND (HALOGEN? OR CHLORINE OR CHLORO? OR BROMINE OR BROMO? L6

2 S ?CHLOROHYPERFORIN OR ?BROMOHYPERFORIN OR ?FLUOROHYPERFORIN OR

7 S ?HYDROHYPERFORIN OR ?HYDROADHYPERFORIN

rs2 S L8 AND (CHLORO OR BROMO OR IODO OR FLUORO OR HALOGEN?) L9

48 S L6 AND (DEPRESSION OR ANTIDEPRESS? OR ALZHEIMER?) L10

26 S L10 AND PHARMACEUTICAL? L11

O S HALOGENATED HYPERFORIN OR HALOGENATED L12

```
chain nodes :
   7 8 9 10 11 12 13 14 15 16 17 18
                                          19
                                              22
                                                  23
                                                      24
                                                         25
                                                             27
                                                                 28
                                                                    29
                       36 37 38 39 40
   30 31 32
             33
                34 35
ring nodes :
                   20
                      21
                         26
   1 2 3
              5
                6
chain bonds :
```

C:\Program Files\Stnexp\Queries\894.str

```
14-15 14-18 15-16 15-17 21-27
                                    21-28
                                           22-23 23-24 23-25 26-29 27-34
   29-30 30-31 31-32 31-33 34-35
                                                 36-38
                                    35-36
                                           36-37
ring bonds :
   1-2 1-6 2-3 2-20 3-4 4-5 5-6 6-21
                                          20-26
                                                 21-26
exact/norm bonds :
   1-19 3-40 5-7 9-10 14-18 15-16 22-23 30-31 35-36
exact bonds :
                                          4-39 5-6 6-14 6-21
                                                               8-9
                      2-20 3-4 4-8 4-5
   1-2 1-6 2-3 2-13
                       15-17 20-26 21-26
   10-12 13-22 14-15
                                           21-27
                                                21-28 23-24 23-25
                       31-33 34-35
                                   36-37
                                           36-38
   27-34
          29-30 31-32
isolated ring systems :
   containing 1 :
G1:H,O
Match level :
   1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS
   10:CLASS 11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS
                      19:CLASS 20:Atom 21:Atom 22:CLASS 23:CLASS
   17:CLASS
            18:CLASS
                      26:Atom 27:CLASS 28:CLASS 29:CLASS 30:CLASS
             25:CLASS
   24:CLASS
                              34:CLASS 35:CLASS 36:CLASS 37:CLASS
   31:CLASS
             32:CLASS
                      33:CLASS
   38:CLASS
           39:CLASS
                      40:CLASS
Element Count :
   Node 16: Limited
       C, C1-3
```

1-19 2-13 3-40 4-8 4-39 5-7 6-14 8-9 9-10 10-11

10-12

STRUCTURE UPLOADED L1

=> d

L1 HAS NO ANSWERS

L1

STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> s 11

SAMPLE SEARCH INITIATED 18:24:17 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED -

100.0% PROCESSED

2 ITERATIONS

2 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:

ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS:

2 TO

PROJECTED ANSWERS:

2 TO

124

L2

2 SEA SSS SAM L1

=> d scan

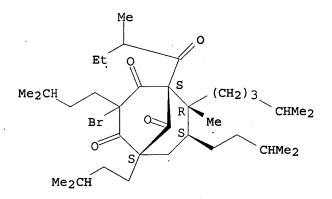
L2 2 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN

Bicyclo[3.3.1] nonane-2,4,9-trione, 3-bromo-6-methyl-1,3,7-tris(3-IN

methylbutyl) -5-(2-methyl-1-oxobutyl) -6-(4-methylpentyl) -, (1S,5S,6R,7S) -(9CI)

C36 H61 Br O4 MF

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L2 REGISTRY COPYRIGHT 2005 ACS on STN

IN Bicyclo [3.3.1] nonane-2,4,9-trione, 3-chloro-6-methyl-1,3,7-tris(3methylbutyl)-5-(2-methyl-1-oxobutyl)-6-(4-methylpentyl)-, (1S,5S,6R,7S)-(9CI)

C36 H61 Cl O4 MF

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> s l1 full FULL SEARCH INITIATED 18:24:33 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 8 TO ITERATE

8 SEA SSS FUL L1

100.0% PROCESSED

8 ITERATIONS

8 ANSWERS

SEARCH TIME: 00.00.01

=> d scan

L3

L3 8 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN

MF C35 H51 Cl O4

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):7

L3 8 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN

MF C36 H61 Cl O4

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 8 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN

MF C35 H51 Br O4

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 8 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN

IN Bicyclo[3.3.1]nonane-2,4,9-trione, 3-bromo-6-methyl-1,3,7-tris(3-methylbutyl)-5-(2-methyl-1-oxobutyl)-6-(4-methylpentyl)-, (1S,5S,6R,7S)-

10/511,894

(9CI)

MF C36 H61 Br O4

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 8 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN

Bicyclo[3.3.1]nonane-2,4,9-trione, 3-chloro-6-methyl-1,3,7-tris(3-methylbutyl)-5-(2-methyl-1-oxopropyl)-6-(4-methylpentyl)-, (1S,5S,6R,7S)(9CI)

MF C35 H59 Cl O4

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 8 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN

MF C36 H53 Cl O4

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 8 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN

MF C35 H59 Br O4

Absolute stereochemistry.

A = A + A + A + A

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

u nomar ∈

L3 8 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN

MF C36 H53 Br O4

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

WEST Search History

Hide Items	Restore	Clear	Cancel

DATE: Monday, December 19, 2005

Hide?	<u>Set</u> <u>Name</u>	Query	<u>Hit</u> <u>Count</u>
	DB=PG	PB; PLUR=YES; OP=ADJ	•
	L2	11 and (halogen or halogenated or chloro or bromo or iodo or fluoro.CLM.)	9
	L1	hyperforin or adhyperforin or \$hydrohyperforin or \$hydroadhyperforin.CLM.	33

END OF SEARCH HISTORY